



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 10:21 AM GMT

PDB ID : 4C00
Title : Crystal structure of TamA from E. coli
Authors : Gruss, F.; Zaehringer, F.; Jakob, R.P.; Burmann, B.M.; Hiller, S.; Maier, T.
Deposited on : 2013-07-30
Resolution : 2.25 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

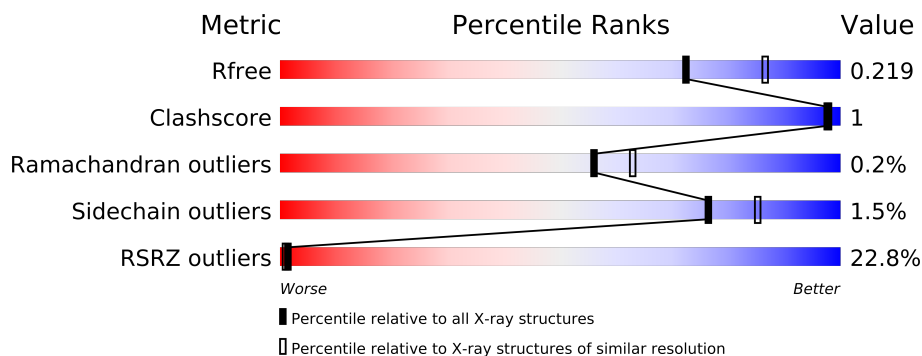
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance


The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1108 (2.28-2.24)
Clashscore	79885	1326 (2.28-2.24)
Ramachandran outliers	78287	1291 (2.28-2.24)
Sidechain outliers	78261	1291 (2.28-2.24)
RSRZ outliers	66119	1110 (2.28-2.24)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	559	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	MC3	A	616	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 9926 atoms, of which 4883 are hydrogens and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

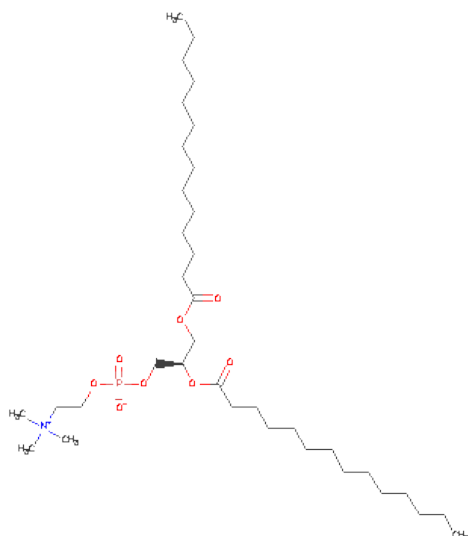
- Molecule 1 is a protein called TRANSLOCATION AND ASSEMBLY MODULE TAMA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	544	8695	2792	4306	760	832	5	0	10	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	GLY	-	EXPRESSION TAG	UNP P0ADE4
A	20	ILE	-	EXPRESSION TAG	UNP P0ADE4
A	21	GLN	-	EXPRESSION TAG	UNP P0ADE4

- Molecule 2 is 1,2-DIMYRISTOYL-RAC-GLYCERO-3-PHOSPHOCHOLINE (three-letter code: MC3) (formula: C₃₆H₇₂NO₈P).



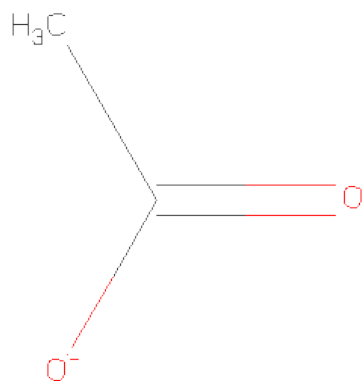
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	118	36	72	1	8	1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C H N O P 118 36 72 1 8 1	0	0
2	A	1	Total C H O P 99 31 59 8 1	0	0
2	A	1	Total C H 30 10 20	0	0
2	A	1	Total C H O P 15 4 5 5 1	0	0
2	A	1	Total C H O P 57 18 31 7 1	0	0
2	A	1	Total C H O P 21 6 7 7 1	0	0
2	A	1	Total C H O 95 31 59 5	0	0
2	A	1	Total C H O 25 9 14 2	0	0
2	A	1	Total C H O P 49 15 25 8 1	0	0
2	A	1	Total C H N O P 107 33 64 1 8 1	0	0
2	A	1	Total C H O P 58 18 31 8 1	0	0
2	A	1	Total C H N O P 71 21 40 1 8 1	0	0
2	A	1	Total C H O P 78 25 44 8 1	0	0
2	A	1	Total C H N O P 24 5 13 1 4 1	0	0
2	A	1	Total C H 23 8 15	0	0

- Molecule 3 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			7	2	3	2		
3	A	1	Total	C	H	O	0	0
			7	2	3	2		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Cl	0	0
			3	3		

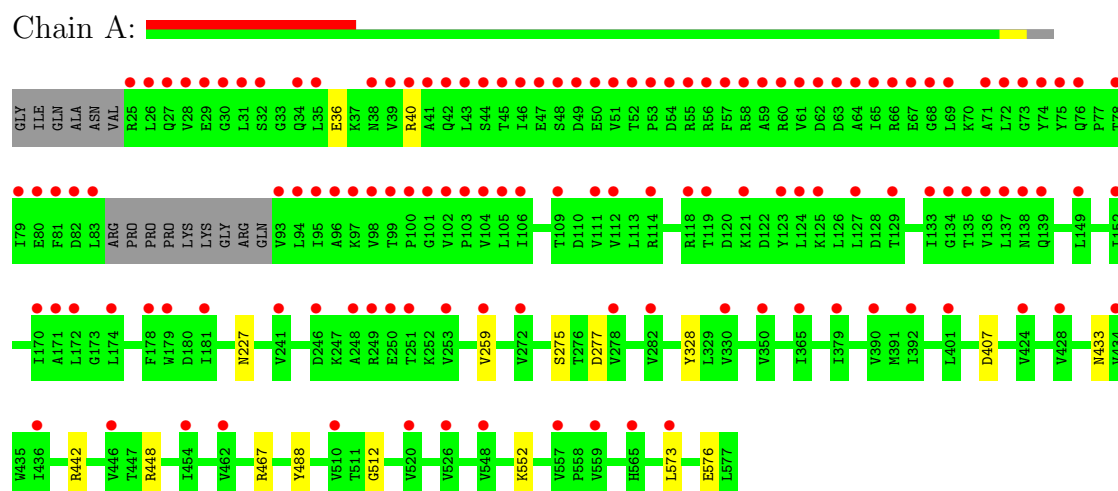
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	226	Total	O	0	0
			226	226		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: TRANSLOCATION AND ASSEMBLY MODULE TAMA



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	77.47Å 261.06Å 57.84Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.27 – 2.25 74.27 – 2.25	Depositor EDS
% Data completeness (in resolution range)	98.4 (74.27-2.25) 98.4 (74.27-2.25)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.25Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.190 , 0.217 0.198 , 0.219	Depositor DCC
R_{free} test set	1946 reflections (3.48%)	DCC
Wilson B-factor (Å ²)	41.1	Xtriage
Anisotropy	0.482	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 60.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 55995 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9926	wwPDB-VP
Average B, all atoms (Å ²)	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, MC3, ACT

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.23	0/4536	0.45	0/6145

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4389	4306	0	7	0
2	A	417	571	2	1	0
3	A	8	6	0	0	0
4	A	3	0	0	2	0
5	A	226	0	0	1	0
All	All	5043	4883	2	7	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 1.

All (7) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:488:TYR:OH	4:A:801:CL:CL	2.32	0.84
1:A:448[B]:ARG:NH2	5:A:2161:HOH:O	2.23	0.70
1:A:328:TYR:OH	2:A:615:MC3:O2P	2.22	0.58
1:A:36:GLU:O	1:A:40:ARG:NH1	2.40	0.55
1:A:275:SER:OG	1:A:277:ASP:OD1	2.29	0.51
1:A:467:ARG:NH1	4:A:801:CL:CL	2.78	0.50
1:A:448[A]:ARG:NH1	1:A:576:GLU:O	2.53	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	551/559 (99%)	525 (95%)	25 (4%)	1 (0%)	56 63

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	512	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	473/475 (100%)	464 (98%)	9 (2%)	69 78

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	227	ASN

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Mol	Chain	Res	Type
1	A	259	VAL
1	A	407	ASP
1	A	433	ASN
1	A	442	ARG
1	A	552[A]	LYS
1	A	552[B]	LYS
1	A	552[C]	LYS
1	A	573	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	139	GLN
1	A	207	GLN
1	A	208	ASN
1	A	227	ASN
1	A	244	GLN
1	A	309	GLN
1	A	322	ASN
1	A	340	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 21 ligands modelled in this entry, 3 are monoatomic - leaving 18 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MC3	A	601	-	45,45,45	1.00	1 (2%)	53,53,53	0.76	2 (3%)
2	MC3	A	602	-	45,45,45	0.97	1 (2%)	53,53,53	0.81	3 (5%)
2	MC3	A	603	-	39,39,45	2.42	1 (2%)	44,44,53	1.96	5 (11%)
2	MC3	A	604	-	8,9,45	2.72	1 (12%)	7,8,53	0.27	0
2	MC3	A	605	-	7,9,45	5.74	1 (14%)	9,12,53	3.95	4 (44%)
2	MC3	A	606	-	24,25,45	3.13	1 (4%)	28,29,53	2.24	6 (21%)
2	MC3	A	607	-	11,13,45	4.57	1 (9%)	14,17,53	3.15	5 (35%)
2	MC3	A	608	-	35,35,45	0.50	1 (2%)	37,37,53	0.85	3 (8%)
2	MC3	A	609	-	8,10,45	2.76	2 (25%)	8,11,53	0.98	1 (12%)
2	MC3	A	610	-	23,23,45	3.77	3 (13%)	28,28,53	2.62	8 (28%)
2	MC3	A	611	-	42,42,45	1.87	3 (7%)	50,50,53	0.96	5 (10%)
2	MC3	A	612	-	26,26,45	3.61	4 (15%)	31,31,53	2.41	8 (25%)
2	MC3	A	613	-	30,30,45	2.21	4 (13%)	38,38,53	0.91	2 (5%)
2	MC3	A	614	-	33,33,45	3.23	3 (9%)	38,38,53	2.08	6 (15%)
2	MC3	A	615	-	10,10,45	3.21	2 (20%)	15,15,53	2.08	2 (13%)
2	MC3	A	616	-	6,7,45	2.93	1 (16%)	5,6,53	0.38	0
3	ACT	A	701	-	1,3,3	1.20	0	0,3,3	0.00	-
3	ACT	A	702	-	1,3,3	1.21	0	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	A	601	-	-	0/49/49/49	0/0/0/0
2	MC3	A	602	-	-	0/49/49/49	0/0/0/0
2	MC3	A	603	-	-	0/41/41/49	0/0/0/0
2	MC3	A	604	-	-	0/7/7/49	0/0/0/0
2	MC3	A	605	-	-	0/8/8/49	0/0/0/0
2	MC3	A	606	-	-	0/26/26/49	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	A	607	-	-	2/13/13/49	0/0/0/0
2	MC3	A	608	-	-	0/37/37/49	0/0/0/0
2	MC3	A	609	-	-	0/9/9/49	0/0/0/0
2	MC3	A	610	-	-	2/25/25/49	0/0/0/0
2	MC3	A	611	-	-	0/46/46/49	0/0/0/0
2	MC3	A	612	-	-	0/28/28/49	0/0/0/0
2	MC3	A	613	-	-	0/34/34/49	0/0/0/0
2	MC3	A	614	-	-	2/35/35/49	0/0/0/0
2	MC3	A	615	-	-	0/8/8/49	0/0/0/0
2	MC3	A	616	-	-	0/5/5/49	0/0/0/0
3	ACT	A	701	-	-	0/0/0/0	0/0/0/0
3	ACT	A	702	-	-	0/0/0/0	0/0/0/0

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	614	MC3	P-O4P	14.93	1.77	1.51
2	A	606	MC3	P-O4P	14.92	1.77	1.51
2	A	610	MC3	P-O4P	14.92	1.77	1.51
2	A	605	MC3	P-O4P	14.87	1.77	1.51
2	A	612	MC3	P-O4P	14.84	1.76	1.51
2	A	607	MC3	P-O4P	14.78	1.76	1.51
2	A	603	MC3	P-O4P	14.68	1.76	1.51
2	A	615	MC3	P-O3P	8.49	1.66	1.51
2	A	604	MC3	C20-C19	-7.66	1.52	1.55
2	A	614	MC3	C21-C20	-7.55	1.52	1.55
2	A	611	MC3	C22-C21	-7.39	1.52	1.55
2	A	613	MC3	C36-C35	-7.31	1.52	1.55
2	A	610	MC3	C35-C34	-7.23	1.52	1.55
2	A	609	MC3	C36-C35	-7.22	1.52	1.55
2	A	616	MC3	C18-C17	-7.15	1.52	1.55
2	A	612	MC3	C18-C17	-7.10	1.52	1.55
2	A	612	MC3	C37-C36	-7.09	1.52	1.55
2	A	614	MC3	C41-C40	-7.03	1.52	1.55
2	A	611	MC3	C43-C42	-7.02	1.52	1.55
2	A	613	MC3	C17-C16	-6.97	1.52	1.55
2	A	610	MC3	C17-C16	-6.25	1.52	1.55
2	A	601	MC3	P-O4P	5.46	1.77	1.60
2	A	613	MC3	P-O4P	5.34	1.77	1.60
2	A	611	MC3	P-O4P	5.24	1.76	1.60
2	A	602	MC3	P-O4P	5.07	1.76	1.60
2	A	615	MC3	P-O4P	4.89	1.76	1.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	609	MC3	O2-C2	-2.75	1.42	1.45
2	A	612	MC3	O3-C11	2.19	1.40	1.33
2	A	613	MC3	P-O3P	2.12	1.66	1.60
2	A	608	MC3	O3-C11	2.02	1.39	1.33

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	603	MC3	O4P-P-O2P	-8.33	94.80	112.88
2	A	612	MC3	O4P-P-O2P	-7.99	95.54	112.88
2	A	614	MC3	O4P-P-O2P	-7.92	95.68	112.88
2	A	610	MC3	O4P-P-O1P	-7.44	95.61	112.73
2	A	607	MC3	O4P-P-O1P	-7.33	95.87	112.73
2	A	606	MC3	O4P-P-O1P	-7.32	95.89	112.73
2	A	605	MC3	O4P-P-O1P	-7.31	95.91	112.73
2	A	603	MC3	O2P-P-O1P	7.10	129.06	112.73
2	A	605	MC3	O2P-P-O1P	7.01	128.85	112.73
2	A	614	MC3	O2P-P-O1P	6.95	128.71	112.73
2	A	606	MC3	O2P-P-O1P	6.95	128.71	112.73
2	A	610	MC3	O2P-P-O1P	6.94	128.68	112.73
2	A	607	MC3	O2P-P-O1P	6.89	128.59	112.73
2	A	612	MC3	O2P-P-O1P	6.86	128.52	112.73
2	A	615	MC3	O2P-P-O1P	6.80	128.36	112.73
2	A	610	MC3	C2-O2-C31	5.29	130.96	117.92
2	A	605	MC3	O2-C2-C1	4.34	113.03	106.74
2	A	610	MC3	O4P-P-O2P	-3.80	104.63	112.88
2	A	606	MC3	O4P-P-O2P	-3.74	104.77	112.88
2	A	615	MC3	O4P-P-O1P	-3.72	95.68	105.86
2	A	605	MC3	O4P-P-O2P	-3.67	104.90	112.88
2	A	610	MC3	O2-C31-C32	3.62	119.48	111.56
2	A	607	MC3	O4P-P-O2P	-3.53	105.21	112.88
2	A	614	MC3	O2-C31-C32	3.53	119.29	111.56
2	A	612	MC3	O4P-P-O1P	-3.48	104.73	112.73
2	A	602	MC3	O2P-P-O1P	3.42	129.05	118.72
2	A	603	MC3	O4P-P-O1P	-3.37	104.98	112.73
2	A	614	MC3	O4P-P-O1P	-3.34	105.05	112.73
2	A	601	MC3	O2P-P-O1P	3.27	128.57	118.72
2	A	611	MC3	O2P-P-O1P	3.25	128.53	118.72
2	A	612	MC3	C2-O2-C31	3.25	125.92	117.92
2	A	614	MC3	C2-O2-C31	3.23	125.89	117.92
2	A	607	MC3	C2-O2-C31	3.11	126.10	116.68
2	A	613	MC3	O2P-P-O1P	3.07	127.97	118.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	603	MC3	C2-O2-C31	3.01	125.35	117.92
2	A	602	MC3	O2P-P-O4P	-2.95	96.01	108.19
2	A	601	MC3	O2P-P-O4P	-2.93	96.07	108.19
2	A	612	MC3	O3-C11-C12	2.90	121.05	111.94
2	A	613	MC3	O2P-P-O4P	-2.90	96.23	108.19
2	A	608	MC3	C3-O3-C11	2.89	125.61	117.13
2	A	610	MC3	O2-C31-O31	-2.89	115.91	123.65
2	A	612	MC3	C3-O3-C11	2.89	125.58	117.13
2	A	611	MC3	O3-C11-O11	-2.71	116.02	123.43
2	A	612	MC3	O3-C11-O11	-2.64	116.23	123.43
2	A	611	MC3	O2P-P-O4P	-2.56	97.61	108.19
2	A	611	MC3	C3-O3-C11	2.47	124.36	117.13
2	A	608	MC3	O3-C11-C12	2.44	119.63	111.94
2	A	614	MC3	O2-C31-O31	-2.44	117.12	123.65
2	A	610	MC3	O3-C3-C2	2.38	115.07	108.83
2	A	606	MC3	O3-C11-O11	-2.36	116.99	123.43
2	A	608	MC3	O3-C11-O11	-2.34	117.03	123.43
2	A	609	MC3	C2-O2-C31	2.32	119.12	115.66
2	A	603	MC3	O2-C2-C1	2.30	116.85	108.40
2	A	607	MC3	O3-C3-C2	2.27	110.67	107.06
2	A	606	MC3	C3-O3-C11	2.26	123.75	117.13
2	A	611	MC3	O3-C11-C12	2.22	118.92	111.94
2	A	610	MC3	O2-C2-C3	2.19	116.46	108.40
2	A	612	MC3	O2-C2-C1	2.14	116.25	108.40
2	A	606	MC3	O3-C11-C12	2.12	118.62	111.94
2	A	602	MC3	O2-C2-C1	2.08	116.04	108.40

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	607	MC3	C2-O2-C31-C32
2	A	614	MC3	C2-O2-C31-C32
2	A	610	MC3	C2-O2-C31-C32
2	A	614	MC3	C2-O2-C31-O31
2	A	607	MC3	C2-O2-C31-O31
2	A	610	MC3	C2-O2-C31-O31

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	544/559 (97%)	1.52	129 (23%) 1 1	31, 55, 196, 254	0

All (129) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	93	VAL	19.6
1	A	28	VAL	18.7
1	A	39	VAL	16.8
1	A	51	VAL	15.8
1	A	26	LEU	15.3
1	A	65	ILE	13.9
1	A	61	VAL	13.1
1	A	27	GLN	12.3
1	A	98	VAL	12.1
1	A	94	LEU	11.7
1	A	57	PHE	11.5
1	A	53	PRO	10.3
1	A	43	LEU	10.2
1	A	46	ILE	9.8
1	A	60	ARG	9.3
1	A	52	THR	9.2
1	A	95	ILE	8.8
1	A	48	SER	8.6
1	A	35	LEU	8.3
1	A	30	GLY	7.7
1	A	96	ALA	7.6
1	A	25	ARG	7.5
1	A	81	PHE	7.4
1	A	83	LEU	7.4
1	A	29	GLU	7.3
1	A	79	ILE	7.2
1	A	54	ASP	7.1

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Mol	Chain	Res	Type	RSRZ
1	A	170	ILE	7.0
1	A	64	ALA	7.0
1	A	106	ILE	6.9
1	A	102	VAL	6.5
1	A	42	GLN	6.4
1	A	97	LYS	6.3
1	A	58	ARG	6.3
1	A	69	LEU	6.2
1	A	59	ALA	6.1
1	A	45	THR	6.1
1	A	31	LEU	6.1
1	A	72	LEU	5.9
1	A	104	VAL	5.8
1	A	73	GLY	5.6
1	A	56	ARG	5.5
1	A	40	ARG	5.3
1	A	49	ASP	5.2
1	A	74	TYR	5.1
1	A	100	PRO	5.1
1	A	136	VAL	5.0
1	A	50	GLU	4.8
1	A	55	ARG	4.7
1	A	38	ASN	4.6
1	A	111	VAL	4.6
1	A	548	VAL	4.5
1	A	112	VAL	4.4
1	A	133	ILE	4.3
1	A	82	ASP	4.2
1	A	137	LEU	4.2
1	A	41	ALA	4.1
1	A	75	TYR	4.1
1	A	152	ILE	3.8
1	A	47	GLU	3.7
1	A	78	THR	3.7
1	A	138	ASN	3.6
1	A	139	GLN	3.6
1	A	124	LEU	3.5
1	A	66	ARG	3.5
1	A	44	SER	3.4
1	A	99	THR	3.4
1	A	71	ALA	3.4
1	A	251	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	103	PRO	3.2
1	A	68	GLY	3.2
1	A	127	LEU	3.2
1	A	129	THR	3.1
1	A	179	TRP	3.1
1	A	171	ALA	3.1
1	A	446	VAL	3.0
1	A	105	LEU	3.0
1	A	34	GLN	3.0
1	A	241	VAL	2.9
1	A	32	SER	2.9
1	A	67	GLU	2.9
1	A	181	ILE	2.9
1	A	125	LYS	2.8
1	A	278	VAL	2.8
1	A	62	ASP	2.8
1	A	253	VAL	2.8
1	A	119	THR	2.7
1	A	436	ILE	2.7
1	A	249	ARG	2.6
1	A	135	THR	2.6
1	A	80	GLU	2.6
1	A	134	GLY	2.6
1	A	250	GLU	2.5
1	A	149	LEU	2.5
1	A	428	VAL	2.5
1	A	462	VAL	2.5
1	A	109	THR	2.5
1	A	272	VAL	2.4
1	A	434	VAL	2.4
1	A	118	ARG	2.4
1	A	401	LEU	2.4
1	A	392	ILE	2.4
1	A	123	TYR	2.3
1	A	573	LEU	2.3
1	A	178	PHE	2.3
1	A	520	VAL	2.3
1	A	248	ALA	2.3
1	A	259	VAL	2.2
1	A	101	GLY	2.2
1	A	559	VAL	2.2
1	A	565	HIS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	390	VAL	2.2
1	A	76	GLN	2.2
1	A	172	LEU	2.2
1	A	365	ILE	2.1
1	A	350	VAL	2.1
1	A	246	ASP	2.1
1	A	454	ILE	2.1
1	A	63	ASP	2.1
1	A	282	VAL	2.1
1	A	424	VAL	2.1
1	A	121	LYS	2.1
1	A	557	VAL	2.1
1	A	379	ILE	2.1
1	A	114	ARG	2.0
1	A	330	VAL	2.0
1	A	510	VAL	2.0
1	A	526	VAL	2.0
1	A	174	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MC3	A	616	8/46	0.19	2.62	62,80,99,99	0
2	MC3	A	602	46/46	0.23	1.74	55,84,141,144	0
2	MC3	A	610	24/46	0.19	1.23	64,112,141,148	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	MC3	A	614	34/46	0.20	1.20	71,104,157,508	0
2	MC3	A	603	40/46	0.21	1.15	41,71,124,507	0
2	MC3	A	601	46/46	0.22	1.11	44,69,102,106	0
2	MC3	A	611	43/46	0.20	1.11	40,74,130,141	0
2	MC3	A	613	31/46	0.20	1.09	44,74,102,115	0
2	MC3	A	604	10/46	0.19	0.49	50,73,98,100	0
2	MC3	A	608	36/46	0.20	0.46	60,85,140,168	0
4	CL	A	803	1/1	0.19	0.29	86,86,86,86	0
2	MC3	A	607	14/46	0.17	0.13	83,121,140,210	0
2	MC3	A	612	27/46	0.18	-0.05	52,102,144,355	0
2	MC3	A	615	11/46	0.16	-0.15	58,94,123,123	0
4	CL	A	802	1/1	0.17	-0.20	51,51,51,51	0
2	MC3	A	609	11/46	0.19	-0.21	51,92,120,129	0
3	ACT	A	702	4/4	0.15	-0.47	49,59,62,67	0
2	MC3	A	606	26/46	0.18	-0.53	48,80,119,129	0
2	MC3	A	605	10/46	0.14	-0.54	91,130,154,154	0
3	ACT	A	701	4/4	0.12	-0.72	54,56,65,65	0
4	CL	A	801	1/1	0.10	-1.05	93,93,93,93	0

6.5 Other polymers ⓘ

There are no such residues in this entry.